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FILE 'REGISTRY' ENTERED AT 15:51:04 ON 07 APR 2011 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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=> D L21 QUE

L6 158 SEA FILE=HCA SPE=ON ABB=ON PLU=ON ("ANTIPAN-LARA, JUAN"/AU OR "GANESHAMURUGAN, SUBRAMANIAM"/AU OR "KATHIRGAMANATHAN,

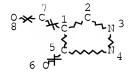
POOPATHY"/AU OR "KUMARAVERL, MUTTULINGHAM"/AU OR "PARAMASWARA, GNANAMOLY"/AU OR "PARTHEEPAN, ARUMUGAM"/AU OR "PRICE, RICHARD"/AU OR "SELVARANJAN, SELVADURAI"/AU OR "SURENDRAKUMAR, SIVAGNANA

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SUNDRAM"/AU)

L7 5017 SEA FILE=HCA SPE=ON ABB=ON PLU=ON ("MERCK PATENT CO

L8 STR



NODE ATTRIBUTES:

DEFAULT MLEVEL IS ATOM

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 8

STEREO ATTRIBUTES: NONE

L9 STR

м 1

NODE ATTRIBUTES:

NSPEC IS RC AT 1 DEFAULT MLEVEL IS ATOM

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 1

STEREO ATTRIBUTES: NONE

L11	3790 SEA FILE=REGISTRY SSS FUL L8 AND L9
L12	151 SEA FILE=REGISTRY SPE=ON ABB=ON PLU=ON L11 AND (RU OR RH OR
	PD OR OS OR PT)/ELS
L13	22 SEA FILE=HCA SPE=ON ABB=ON PLU=ON L12
L14	32 SEA FILE=REGISTRY SPE=ON ABB=ON PLU=ON L11 AND IR/ELS
L15	27 SEA FILE=REGISTRY SPE=ON ABB=ON PLU=ON L14 NOT S/ELS

### => FIL HCA

FILE 'HCA' ENTERED AT 15:51:25 ON 07 APR 2011
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# => D L18 1-4 IBIB ABS HITSTR HITIND RETABLE

L18 ANSWER 1 OF 4 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 145:198513 HCA Full-text

TITLE: Electroluminescent device fabrication by spin coating

electroluminescent organometallic complexes on coated

substrates

INVENTOR(S): Kathirgamanathan, Poopathy;

Ganeshamurugan, Subramaniam; Price,

Richard

PATENT ASSIGNEE(S): Oled-T Limited, UK SOURCE: PCT Int. Appl., 51 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.					KIN	D	DATE			APPLICATION NO.				DATE			
WO	2006	0774	02		A1	_									2	0060	 119
	W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	ΚE,	KG,	KM,	KN,	ΚP,	KR,
		ΚZ,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	LY,	MA,	MD,	MG,	MK,	MN,	MW,	MX,
		MZ,	NA,	NG,	NΙ,	NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,
		SG,	SK,	SL,	SM,	SY,	ΤJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC,
		VN,	YU,	ZA,	ZM,	ZW											
	RW:	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FΙ,	FR,	GB,	GR,	HU,	IE,
		IS,	IT,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	BJ,
		CF,	CG,	CI,	CM,	GA,	GN,	GQ,	GW,	ML,	MR,	ΝE,	SN,	TD,	TG,	BW,	GH,
		GM,	ΚE,	LS,	MW,	MΖ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,	AZ,	BY,
		KG,	KΖ,	MD,	RU,	ТJ,	TM										
EP	1839	464			A1		2007	1003		EP 2	006-	7027	71		2	0060	119
	R:	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FΙ,	FR,	GB,	GR,	HU,	IE,
		IS,	ΙΤ,	LI,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR	
	1011															0060	119
JP	2008	5292	12		T		2008	0731		JP 2	007-	5517.	36		2	0060	119
US	2008	0160	182		A1		2008	0703		US 2	007-	7950	07		2	0070	710

IN 2007DN05397 Α 20070817 IN 2007-DN5397 20070712 KR 2007102556 Α 20071018 KR 2007-7018852 20070817 PRIORITY APPLN. INFO.: GB 2005-1426 A 20050122 W WO 2006-GB169 20060119

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): MARPAT 145:198513

AB Methods of forming electroluminescent devices are described which entail depositing by spin coating a layer of an electroluminescent organometallic complex on a substrate (which is the anode) which is coated with a layer of a polymer. The polymer is preferably a conductive or charge-transporting polymer or material.

IT 647838-95-7

(electroluminescent device fabrication by spin coating electroluminescent organometallic complexes on coated substrates)

RN 647838-95-7 HCA

CN Iridium,  $[4-[3,3-dimethyl-1-(oxo-\kappa0)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa03$ ]bis $[2-(2-pyridinyl-\kappa N)phenyl-\kappa0]-$  (CA INDEX NAME)

IPCI H05B0033-14 [I,A]; H01L0051-50 [I,A]; H01L0051-56 [I,A]; C09K0011-06
[I,A]; H01L0051-30 [I,A]

IPCR H05B0033-14 [I,A]; C09K0011-06 [I,C]; C09K0011-06 [I,A]; H01L0051-00
 [I,C\*]; H01L0051-00 [I,A]; H01L0051-05 [I,C]; H01L0051-30 [I,A];
 H01L0051-50 [I,C]; H01L0051-50 [I,A]; H01L0051-56 [I,A]; H05B0033-14 [I,C]

CC 73-11 (Optical, Electron, and Mass Spectroscopy and Other Related Properties)

Section cross-reference(s): 76

86-73-7D, 9H-Fluorene, derivs. 159-66-0D, 9,9'-Spirobi[9H-fluorene], ΤT derivs. 193-44-2 905-62-4 1217-45-4, 9,10-Dicyanoanthracene 2085-33-8, Tris(8-hydroxyquinolinato)aluminum 4733-39-5, Bathocuproin 5521-31-3D, derivs. 7429-90-5, Aluminum, uses 7439-93-2, Lithium, uses 7439-95-4, Magnesium, uses 7440-03-1D, Niobium, compds. 7440-04-2D, Osmium, compds. 7440-05-3D, Palladium, compds. 7440-06-4D, Platinum, 7440-16-6D, Rhodium, compds. 7440-18-8D, Ruthenium, compds. 7440-25-7D, Tantalum, compds. 7440-32-6D, Titanium, compds. 7440-39-3, Barium, uses 7440-58-6D, Hafnium, compds. 7440-62-2D, Vanadium, 7440-70-2, Calcium, uses 7789-24-4, Lithium fluoride, uses compds. 17595-05-0 19414-67-6 23467-27-8 25067-59-8, 15082-28-7 Poly(vinylcarbazole) 25135-15-3D, derivs. 25233-30-1, Polyaniline 25387-93-3 26009-24-5, Poly(p-phenylenevinylene) - 31366-25-3D, derivs. 37271-44-6 58280-31-2 58328-31-7, CBP 58328-31-7D, derivs. 65181-78-4, N,N'-Diphenyl-N,N'-bis(3-methylphenyl)-1,1'-biphenyl-4,4'diamine 66946-48-3D, derivs. 95270-88-5D, derivs. 98038-22-3, Aniline-m-sulfanilic acid copolymer 121220-44-8, 124729-98-2 o-Ethylaniline-o-toluidine copolymer 123847-85-8

126415-16-5, Aniline-o-anisidine copolymer 126415-18-7, o-Aminophenol-aniline copolymer 126415-20-1, o-Aminophenol-o-toluidine copolymer 126415-22-3, o-Phenylenediamine-o-toluidine copolymer 135804-06-7 138372-67-5 142289-08-5D, derivs. 146162-54-1 148044-16-0 148896-39-3 150405-69-9 157755-87-8 203642-12-0D, derivs. 214341-85-2D, derivs. 221455-80-7 300576-41-4 432042-07-4 432042-08-5 474974-61-3 474974-62-4 647838-95-7 861532-86-7D, [9,9'-Bianthracene]-10,10'-diamine, N-aryl derivs. 863714-50-5 902119-35-1

(electroluminescent device fabrication by spin coating electroluminescent organometallic complexes on coated substrates)

#### RETABLE

Referenced Author	Year   VOL		Referenced Work	Referenced
(RAU)	(RPY) (RVL) 		(RWK) :+==========	File
	-+ <del>-</del>	-т		т
Eriyama, Y	2004	1	US 2004106006 A1	HCA
Kamatant, J	[2003 ]	1	US 2003224208 A1	HCA
Kathirgamanathan, P	2004	1	WO 2004084325 A	HCA

L18 ANSWER 2 OF 4 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 144:458233 HCA Full-text

TITLE: Electroluminescent devices with anode buffer layers

INVENTOR(S): Kathirgamanathan, Poopathy;

Ganeshamurugan, Subramaniam; Kumaraverl,

Muttulingham; Partheepan, Arumugam;

Paramaswara, Gnanamoly Nuko 70 Limited, UK

PATENT ASSIGNEE(S): Nuko 70 Limited, UK SOURCE: PCT Int. Appl., 89 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

E	PAT	ENT I	. OI			KIN	D :				APPL	ICAT	ION I	NO.		D	ATE	
V	 VO	2006	0486	 35		A1				1	WO 2	005-	GB42:	 22		2	0051	101
		W:	ΑE,	AG,	AL,	AM,	AT,	AU,	AZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,
			CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,
			GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KΕ,	KG,	KM,	KN,	KP,	KR,
			KΖ,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	LY,	MA,	MD,	MG,	MK,	MN,	MW,	MX,
			MZ,	ΝA,	NG,	NI,	NO,	NΖ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,
			SG,	SK,	SL,	SM,	SY,	ТJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC,
			VN,	YU,	ZA,	ZM,	ZW											
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			IS,	ΙT,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR,	BF,	ВJ,
			CF,	CG,	CI,	CM,	GA,	GN,	GQ,	GW,	ML,	MR,	NE,	SN,	TD,	TG,	BW,	GH,
			GM,	ΚE,	LS,	MW,	MZ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,	AZ,	BY,
			KG,	KΖ,	MD,	RU,	ТJ,	TM										
E	ΞP	1812	530			A1		2007	0801		EP 2	005-	8001	28		2	0051	101
		R:	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,	FI,	FR,	GB,	GR,	HU,	ΙE,
			IS,	ΙT,	LI,	LT,	LU,	LV,	MC,	NL,	PL,	PT,	RO,	SE,	SI,	SK,	TR	
-	JΡ	2008	5194	27		T		2008	0605		JP 2	007-	5385	21		2	0051	101
J	JS	2008	0199	727		A1		2008	0821	1	US 2	007-	6667	66		2	0070	625
PRIOR]	ΙΤΥ	APP	LN.	INFO	.:						GB 2	004 -	2429	4	Ž	A 2	0041	103
										1	WO 2	005-	GB 42	22	1	w 2	0051	101

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

AB Electroluminescent devices are described which are provided with a buffer layer on the anode, the buffer material being selected from metal tetra-p-tolyl porphinato complexes and bianthryl compds. [9,9'-Bianthracene]-10,10'-

diamine, N,N'-di-2-naphthalenyl-N,N'-diphenyl- [223735-42-0] or [9,9'-Bianthracene]-10,10'-diamine, N,N'-di-1-naphthalenyl-N,N'-diphenyl-. The electroluminescent materials may be organometallic compds., including multinuclear complexes.

IT 647838-95-7

(electroluminescent devices with anode buffer layers)

RN 647838-95-7 HCA

CN Iridium, [4-[3,3-dimethyl-1-(oxo-κ0)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-κ03]bis[2-(2-pyridinyl-κN)phenyl-κC]- (CA INDEX NAME)

IPCI C09K0011-06 [I,A]; H01L0051-50 [I,A]
IPCR C09K0011-06 [I,A]; C09K0011-06 [I,C]; H01L0051-00 [I,C\*]; H01L0051-00
[I,A]; H01L0051-50 [I,C]; H01L0051-50 [I,A]; H05B0033-14 [I,C\*];
H05B0033-14 [I,A]; H05B0033-22 [I,C\*]; H05B0033-22 [I,A]

CC 73-11 (Optical, Electron, and Mass Spectroscopy and Other Related Properties)

Section cross-reference(s): 76

ΙT 2085-33-8, Tris(8-hydroxyquinolinato)aluminum 4733-39-5, Bathocuproin 7429-90-5, Aluminum, uses 7429-90-5D, Aluminum, compds. 7429-91-6D, Dysprosium, compds. 7439-88-5D, Iridium, compds. 7439-89-6D, Iron, 7439-92-1D, Lead, compds. 7439-93-2, Lithium, uses compds. 7439-93-2D, Lithium, compds. 7439-94-3D, Lutetium, compds. 7439-95-4, Magnesium, uses 7439-95-4D, Magnesium, compds. 7439-96-5D, Manganese, 7439-98-7D, Molybdenum, compds. 7440-00-8D, Neodymium, compds. 7440-02-0D, Nickel, compds. 7440-03-1D, Niobium, compds. 7440-04-2D, Osmium, compds. 7440-05-3D, Palladium, compds. 7440-06-4D, Platinum, 7440-09-7D, Potassium, compds. 7440-10-0D, Praseodymium, 7440-12-2D, Promethium, compds. 7440-16-6D, Rhodium, compds. 7440-17-7D, Rubidium, compds. 7440-18-8D, Ruthenium, compds. 7440-19-9D, Samarium, compds. 7440-22-4D, Silver, compds. 7440-23-5D, Sodium, compds. 7440-24-6D, Strontium, compds. 7440-25-7D, Tantalum, 7440-27-9D, Terbium, compds. 7440-30-4D, Thulium, compds. 7440-31-5D, Tin, compds. 7440-32-6D, Titanium, compds. 7440-39-3D, Barium, compds. Antimony, compds. 7440-41-7D, Beryllium, 7440-43-9D, Cadmium, compds. 7440-42-8D, Boron, compds. 7440-45-1D, Cerium, compds. 7440-46-2D, Cesium, compds. 7440-47-3D, 7440-48-4D, Cobalt, compds. Chromium, compds. 7440-50-8D, Copper, 7440-52-0D, Erbium, compds. 7440-53-1D, Europium, compds. compds. 7440-54-2D, Gadolinium, compds. 7440-55-3D, Gallium, compds. 7440-56-4D, Germanium, compds. 7440-57-5D, Gold, compds. Holmium, compds. 7440-61-1D, Uranium, compds. 7440-62-2D, Vanadium, 7440-64-4D, Ytterbium, compds. 7440-65-5D, Yttrium, compds. 7440-67-7D, Zirconium, compds. 7440-70-2, Calcium, uses 7440-70-2D, Calcium, compds. 7440-74-6D, Indium, compds. 14405-36-8 15133-54-7

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17457-88-4 21333-45-9 21392-78-9 23467-27-8 24911-10-2 25067-59-8, Poly(vinylcarbazole) 26009-24-5, Poly(p-phenylenevinylene) 37271-44-6 50851-57-5 58280-31-2 58328-31-7, CBP 63448-47-5 65181-78-4, N,N'-Diphenyl-N,N'-bis(3-methylphenyl)-1,1'-biphenyl-4,4'-diamine 135804-06-7 142289-08-5 146162-54-1 148896-39-3 223735-42-0 386223-21-8 432042-07-4 $47838-95-7 873428-96-7 885502-27-2 885502-29-4 885502-30-7 885502-32-9 (electroluminescent devices with anode buffer layers)
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RETABLE

Referenced Author	·	Referenced Work	Referenced
(RAU)	(RPY) (RVL) (RPG)	(RWK)	File
	=+====+=====	=+===========	=+=======
Anon	1999  1999	PATENT ABSTRACTS OF	1
Aziz, H	[2004 ]	US 2004018381 A1	1
Eastman Kodak Company	1988	EP 0278758 A	HCA
Hu, N	[2003 ]	US 6670054 B1	HCA
Lg Electronics Inc	[2003 ]	EP 1317005 A	HCA
Toyo Ink Mfg Co Ltd	1999	JP 11265788 A	HCA
OS.CITING REF COUNT:	1 THERE ARE 1	1 CAPLUS RECORDS THAT	CITE THIS RECORD
	(1 CITINGS)	)	

L18 ANSWER 3 OF 4 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 143:275247 HCA Full-text

TITLE: Electroluminescent organometallic materials and their

preparation and devices using them
Kathirgamanathan, Poopathy; Price,
Richard; Ganeshamurugan, Subramaniam;
Paramaswara, Gnanamoly; Kumaraverl,
Muttulingham; Partheepan, Arumugam;
Selvaranjan, Selvadurai; Antipan-Lara,

Juan; Surendrakumar, Sivagnanasundram

PATENT ASSIGNEE(S): Elam-T Limited, UK SOURCE: PCT Int. Appl., 66 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

INVENTOR(S):

PATENT NO.	KIND DATE	APPLICATION NO.	DATE
=	A2 20050901 A3 20051103	WO 2005-GB446	20050210
W: AE, AG, A	AL, AM, AT, AU, AZ,	BA, BB, BG, BR, BW,	BY, BZ, CA, CH,
CN, CO, C	CR, CU, CZ, DE, DK,	DM, DZ, EC, EE, EG,	ES, FI, GB, GD,
GE, GH, G	GM, HR, HU, ID, IL,	IN, IS, JP, KE, KG,	KP, KR, KZ, LC,
LK, LR, I	LS, LT, LU, LV, MA,	MD, MG, MK, MN, MW,	MX, MZ, NA, NI,
NO, NZ, C	OM, PG, PH, PL, PT,	RO, RU, SC, SD, SE,	SG, SK, SL, SY,
TJ, TM, T	TN, TR, TT, TZ, UA,	UG, US, UZ, VC, VN,	YU, ZA, ZM, ZW
RW: BW, GH, G	GM, KE, LS, MW, MZ,	NA, SD, SL, SZ, TZ,	UG, ZM, ZW, AM,
AZ, BY, K	KG, KZ, MD, RU, TJ,	TM, AT, BE, BG, CH,	CY, CZ, DE, DK,
EE, ES, F	FI, FR, GB, GR, HU,	IE, IS, IT, LT, LU,	MC, NL, PL, PT,
RO, SE, S	SI, SK, TR, BF, BJ,	CF, CG, CI, CM, GA,	GN, GQ, GW, ML,
MR, NE, S	SN, TD, TG		
EP 1723213	A2 20061122	EP 2005-708271	20050210
R: AT, BE, E	BG, CH, CY, CZ, DE,	DK, EE, ES, FI, FR,	GB, GR, HU, IE,
		PL, PT, RO, SE, SI,	SK, TR
JP 2007524680	T 20070830	JP 2006-552679	20050210
KR 2007004719	A 20070109	KR 2006-7018827	20060914
US 200900090 <b>6</b> 0	A1 20090108	US 2007-589183	20070808

PRIORITY APPLN. INFO.:

GB 2004-3322 A 20040214 WO 2005-GB446 W 20050210 7

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT OTHER SOURCE(S): MARPAT 143:275247

AB Electroluminescent compds. are described by the general formula I, II, and III (R1-6 = independently selected H, (un)substituted hydrocarbyl groups such as (un)substituted aliphatic groups, (un)substituted aromatic, heterocyclic and polycyclic ring structures, fluorocarbons such as trifluoryl Me groups, halogens such as F, or thiophenyl groups; R1, R2 and R3 can form (un)substituted fused aromatic, heterocyclic and polycyclic ring structures and can be copolymerizable with a monomer, e.g. styrene; M = ruthenium, rhodium, palladium, osmium, iridium, or platinum; and n+2 is the valency of M). Methods of preparing the compds. are also described which entail reacting a bridged complex with an appropriate ligand. Electroluminescent devices employing the materials are also described.

IT 647838-95-7P 863714-47-0P 863714-48-1P 863714-49-2P

(electroluminescent organometallic materials and their preparation and devices using them)

RN 647838-95-7 HCA

CN Iridium,  $[4-[3,3-dimethyl-1-(oxo-\kappa0)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa03$ ]bis $[2-(2-pyridinyl-\kappa N)phenyl-\kappa0]-$  (CA INDEX NAME)

RN 863714-47-0 HCA

CN Iridium,  $[4-[3,3-dimethyl-1-(oxo-\kappa O)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa O3$ ]bis[3-fluoro-2-(2-pyridinyl- $\kappa N$ )phenyl- $\kappa C$ ]- (CA INDEX NAME)

RN 863714-48-1 HCA

CN Iridium, bis[3,5-difluoro-2-(2-pyridinyl- $\kappa$ N)phenyl- $\kappa$ C][4-[3,3-dimethyl-1-(oxo- $\kappa$ O)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ O3]- (CA INDEX NAME)

RN 863714-49-2 HCA

CN Iridium, bis[3,5-difluoro-2-(2-pyridinyl-κN)phenyl-κC][4-[3,3-dimethyl-1-(oxo-κO)butyl]-2-(4-fluorophenyl)-2,4-dihydro-5-methyl-3H-pyrazol-3-onato-κO3]- (CA INDEX NAME)

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IPCI C09K0011-06 [ICM,7]; H01L0051-30 [ICS,7]; H01L0051-05 [ICS,7,C*]
IPCR C09K0011-06 [I,C*]; C09K0011-06 [I,A]; H01L0051-00 [N,C*]; H01L0051-00
      [N,A]; H01L0051-05 [I,C*]; H01L0051-30 [I,A]; H01L0051-50 [I,C*];
      H01L0051-50 [I,A]
```

CC 73-5 (Optical, Electron, and Mass Spectroscopy and Other Related Properties)

Section cross-reference(s): 29, 76

IT 7440-04-2DP, Osmium, compds. 7440-05-3DP, Palladium, compds. 7440-06-4DP, Platinum, compds. 7440-16-6DP, Rhodium, compds.

7440-18-8DP, Ruthenium, compds. 647838-95-72

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863714-47-0P 863714-48-1P 863714-49-2P

863714-50-5P 863714-51-6P 863714-52-7P 863714-54-9P

(electroluminescent organometallic materials and their preparation and devices using them)

RETABLE

L18 ANSWER 4 OF 4 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER:  $140:114240 \quad \text{HCA} \quad \underline{\text{Full-text}}$ 

TITLE: Metal chelates in a photovoltaic device

INVENTOR(S): Kathirgamanathan, Poopathy;
Antipan-Lara, Juan; Partheepan,

Arumuqam

PATENT ASSIGNEE(S): Elam-Limited, UK

SOURCE: PCT Int. Appl., 59 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PATENT NO.			KIN	D	DATE APPLICATION NO.					DATE							
		2004				A2 A3		2004		1	WO 2	003-	GB303	35		2	0030	714
	WU	₩:		-		_		AU,		BA,	BB,	BG,	BR,	BY,	BZ,	CA,	CH,	CN,
					•			DK, IN,		•		•	•					
			LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NO,	NZ,	OM,	PH,
			,	,	•	•	•	SE, ZA,	,	•	SL,	TJ,	TM,	TN,	TR,	TT,	TZ,	UA,
		RW:			•			MZ,		•		•						
					•	•	•	TM, IE,	•	•	•	•	•	•	•	•	•	•
			BF,	ΒJ,	CF,	CG,	CI,	CM,	GΑ,	GN,	GQ,	GW,	${ m ML}$ ,	MR,	NE,	SN,	TD,	ΤG
,	ΑU	2003	2810	03		A1		2004	0202		AU 2	003-	2810	03		2	0030	714
PRIOR	PRIORITY APPLN. INFO.:							(	GB 2	002-	1615	4	2	A 2	A 20020712			
										1	WO 2	003-	GB30:	35	I	₩ 2	0030	714

# OTHER SOURCE(S): MARPAT 140:114240

AB A photovoltaic device uses a metal chelate as the photovoltaic element. The device comprises sequentially (1) a first electrode comprising a metal, (2) the photovoltaic element, and (3) a second electrode. The photovoltaic element comprises an organometallic complex with an organic ligand and a metal (a rare earth, transition metal, lanthanide, or an actinide).

IT 647838-95-7

(metal chelates in photovoltaic device)

RN 647838-95-7 HCA

CN Iridium,  $[4-[3,3-dimethyl-1-(oxo-\kappa0)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa$ O3]bis $[2-(2-pyridinyl-\kappa N)phenyl-\kappa C]-$  (CA INDEX NAME)

IPCI H01L0051-20 [ICM,7]; H01L0051-30 [ICS,7]; H01L0051-05 [ICS,7,C\*]
IPCR H01L0051-00 [I,C\*]; H01L0051-00 [I,A]; H01L0051-05 [I,C\*]; H01L0051-30
 [I,A]; H01L0051-40 [N,A]

CC 52-2 (Electrochemical, Radiational, and Thermal Energy Technology) Section cross-reference(s): 74, 76

147-14-8, Copper phthalocyanine 2085-33-8 7429-90-5, Aluminum, uses TΤ 7429-90-5D, Aluminum, complex 7439-88-5D, Iridium, complex 7439-89-6D, Iron, complex 7439-92-1D, Lead, complex 7439-93-2D, Lithium, complex 7439-95-4D, Magnesium, complex 7439-96-5D, Manganese, complex 7439-98-7D, Molybdenum, complex 7440-02-0D, Nickel, complex 7440-03-1D, Niobium, complex 7440-04-2D, Osmium, complex 7440-05-3D, Palladium, complex 7440-06-4D, Platinum, complex 7440-09-7D, Potassium, complex 7440-16-6D, Rhodium, complex 7440-17-7D, Rubidium, 7440-18-8D, Ruthenium, complex 7440-20-2D, Scandium, complex 7440-22-4D, Silver, complex 7440-23-5D, Sodium, complex 7440-24-6D, Strontium, complex 7440-25-7D, Tantalum, complex 7440-31-5D, Tin, complex 7440-32-6D, Titanium, complex 7440-36-0D, Antimony, complex 7440-39-3D, Barium, complex 7440-41-7D, Beryllium, complex 7440-42-8D, Boron, complex 7440-43-9D, Cadmium, complex 7440-46-2D, Cesium, 7440-47-3D, Chromium, complex 7440-48-4D, Cobalt, complex complex 7440-50-8D, Copper, complex 7440-55-3D, Gallium, complex 7440-56-4D, Germanium, complex 7440-57-5D, Gold, complex 7440-62-2D, Vanadium, 7440-65-5D, Yttrium, complex 7440-66-6D, Zinc, complex 7440-67-7D, Zirconium, complex 7440-70-2D, Calcium, complex 7440-74-6D, Indium, complex 7789-24-4, Lithium fluoride, uses 14913-52-1D, Neodymium(3+), complex, uses 15956-38-4 16910-54-6D, Europium (2+), complex, uses 17457-88-4 18472-30-5D, Erbium (3+), complex, uses 18581-58-3 18923-26-7D, Cerium(3+), complex, uses 18923-27-8D, Ytterbium(3+), complex, uses 21392-78-9 22541-14-6D, Praseodymium(3+), complex, uses 22541-16-8D, Promethium(3+), complex, uses 22541-17-9D, Samarium(3+), complex, uses 22541-18-0D, Europium(3+), complex, uses 22541-19-1D, Gadolinium(3+), complex, uses 22541-20-4D, Terbium(3+), complex, uses 22541-21-5D, Dysprosium(3+), complex, uses 22541-22-6D, Holmium(3+), complex, uses 22541-23-7D, Thulium(3+), complex, uses 22541-24-8D, Lutetium(3+), complex, uses 22578-81-0D, Uranium(3+), complex, uses 25387-93-3 65181-78-4, N, N'-Diphenyl-N, N'-bis(3-methylphenyl)-1,1'-biphenyl-4,4'-diamine 114206-51-8 156882-92-7 647838-95-7

(metal chelates in photovoltaic device)

### RETABLE

Referenced Author (RAU)	(RPY)	(RVL)   (RPG)	Referenced Work   (RWK)	Referenced
Anon	_		IWO 0243444 A2	HCA
Anon	i	i	IEP 0556005 A1	HCA
Anon	i	i	US 6153824 A	HCA
Anon	1	1	US 6310282 B1	HCA
OS.CITING REF COUNT:	4	THERE ARE	4 CAPLUS RECORDS THAT	CITE THIS RECORD
		(4 CITINGS)	)	

## => D L20 1-16 IBIB ABS HITSTR HITIND RETABLE

L20 ANSWER 1 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 142:488306 HCA Full-text

TITLE: Association Behavior of 4-Acylpyrazolone Derivative and Tertiary Amine of High Molecular Weight in Antagonistic Synergistic Extraction of Palladium

AUTHOR(S): Zhang, Anyun; Wanyan, Guanghui; Kumagai, Mikio
CORPORATE SOURCE: Nuclear Chemistry and Chemical Engineering Center,
Institute of Research and Innovation (IRI), Kashiwa,

Chiba-ken, 277-0861, Japan

SOURCE: Journal of Solution Chemistry (2004), 33(8),

1017-1028

CODEN: JSLCAG; ISSN: 0095-9782

PUBLISHER: Springer Science+Business Media, Inc.

DOCUMENT TYPE: Journal LANGUAGE: English

To find an effective extraction and removal method for palladium(II), which is one of the main fission products from an acidic nuclear spent fuel solution, the extraction behavior of palladium(II) from a nitric acid medium by an acidic chelating extractant, 1-phenyl-3-methyl-4-trifluoroacetylpyrazolone-5-one (HPMTP) and a tertiary amine of high mol. weight, tri-n-octylamine (TOA), has been studied by spectrophotometry. A noticeable antagonistic extraction effect was observed in the extraction system under the given conditions. To understand this phenomenon, a preliminary investigation was performed to explain the mechanism of this reaction. According to the theory of corresponding solns. (TCS), the association reaction between HPMTP and TOA is proposed in the organic phase. An associated species, HPMTP·TOA, formed through hydrogen bonding in a chloroform medium might be the main reason why an antagonistic extraction effect occurred. The association constant between HPMTP and TOA was calculated to be 2.86 ± 0.05.

IT 851853-56-0P

(preparation of palladium complex with fluoroacylpyrazolone derivative)

RN 851853-56-0 HCA

CN Palladium(2+), bis[2,4-dihydro-5-methyl-2-phenyl-4-(trifluoroacetyl-κO)-3H-pyrazol-3-one-κO3]- (9CI) (CA INDEX NAME)

CC 68-2 (Phase Equilibriums, Chemical Equilibriums, and Solutions) Section cross-reference(s): 71, 73

IT 851853-56-0P

(preparation of palladium complex with fluoroacylpyrazolone derivative)  $\ensuremath{\mathsf{RETABLE}}$ 

Referenced Author (RAU)	Year	Referenced Work   Referenced   (RWK)   File
	=+====+====	=+===============
Horwitz, E	1997  33  25	Reactive and Functio HCA
Horwitz, E	1991  9  1	Solvent Extr Ion Exc HCA
Jensen, B	1959  13  1668	Acta Chem Scand   HCA
Kolarik, Z	1999  17  1155	Solvent Extr Ion Exc HCA
Koma, Y	1998  16  1357	Solvent Extr Ion Exc HCA
Law, J	2001  19  23	Solvent Extr Ion Exc HCA
Law, J	1999  19  27	Waste Management   HCA
Li, X	1993  12  2021	Polyhedron   HCA
Madic, C	2002    27	JAERI-Conf 2002-004   HCA

Margenk, Z	1983	1	1353	Spectrophotometric D	1
Mirza, M	1979	41	772	J Inorg Nucl Chem	HCA
Mukai, H	1997	13	145	Anal Sci	HCA
Nagasaki, S	1994	12	459	Solvent Extr Ion Exc	HCA
Pokhitonov, V	12003	1	1	Proceedings of 2003	1
Reddy, M	12000	18	1135	Solvent Extr Ion Exc	HCA
Romanovskiy, V	2001	19	1	Solvent Extr Ion Exc	HCA
Schulz, W	1988	123	1191	Sep Sci Technol	HCA
Uchiyama, G	12002	1	197	JAERI-Conf 2002-004	HCA
Uehara, A	2001	117	1045	Anal Sci	1
Umetani, S	1985	1	55	Anal Sci	HCA
Wei, Y	12004	41	315	J Nucl Sci Technol	HCA
Wood, D	11997	132	1241	Sep Sci Technol	HCA
Zhang, A	1996		142	Chemistry	HCA
Zhang, A	1996	1	150	Chemistry	1
Zhang, A		172	I	Radiat Phys Chem (in	.1
Zhang, A	12003	21	591	Solvent Extr Ion Exc	HCA
OS.CITING REF COUNT:	8	THE	RE ARE 8	8 CAPLUS RECORDS THAT	CITE THIS RECORD
		(8 (	CITINGS	)	

L20 ANSWER 2 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 141:198940 HCA Full-text

TITLE: A 4-acyl-5-pyrazolone ligand (HQ) in N-unidentate

coordination mode in a Rh(CO)2Cl(HQ)-type complex

AUTHOR(S): Cingolani, Augusto; Marchetti, Fabio; Pettinari,

Claudio; Pettinari, Riccardo; Skelton, Brian W.;

White, Allan H.

CORPORATE SOURCE: Dipartimento di Scienze Chimiche, Universita degli

Studi di Camerino, Camerino, MC, I-62032, Italy

SOURCE: Inorganic Chemistry Communications (2004),

7(2), 235-237

CODEN: ICCOFP; ISSN: 1387-7003

PUBLISHER: Elsevier Science B.V.

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 141:198940

AB Rh(CO) 2Cl(HQ) (HQ = 4-acylpyrazolone), the ligand being bonded in N-unidentate fashion, was synthesized and structurally characterized. [Rh(CO) 2Q] and [CpRhQCl] were also prepared The crystal structure of HQ was also determined

IT 737787-67-6P

(preparation and crystal structure of)

RN 737787-67-6 HCA

CN Rhodium, dicarbonylchloro[2-furanyl(5-hydroxy-3-methyl-1-phenyl-1H-pyrazol- $4-yl-\kappa N2$ )methanone]-, (SP-4-3)- (CA INDEX NAME)

$$\begin{array}{c|c}
O & OH \\
\hline
N & Ph \\
\hline
N & C & Rh + C & O \\
\hline
C1 - & O
\end{array}$$

RN 737787-66-5 HCA

CN Rhodium, dicarbonyl[ $4-(2-furanylcarbonyl-\kappa O)-2$ ,  $4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa O3$ ]-, (SP-4-3)- (CA INDEX NAME)

RN 737787-68-7 HCA

CN Rhodium, chloro[ $4-(2-furanylcarbonyl-\kappa O)-2$ ,  $4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3$ ][ $(1,2,3,4,5-\eta)-1$ ,2,3,4,5-pentamethyl-2,4-cyclopentadien-1-yl]- (CA INDEX NAME)

CC 78-7 (Inorganic Chemicals and Reactions)

Section cross-reference(s): 75

IT 737787-67-6P

(preparation and crystal structure of)

IT 737787-66-5P 737787-68-7P

(preparation of)

RETABLE

Referenced Author (RAU)	(RPY) (RVL)		Referenced Work   (RWK) -+	Referenced   File
Arndtsen, B	1995  270	1970	Science	HCA
Barbera, J	1999  38	13085	Inorg Chem	HCA
Bonati, F	1964	3156	J Chem Soc	HCA
Bonati, F	1985  4	1357	Polyhedron	HCA
Chaloner, P	1994	1	Homogeneous Hydrogen	
Cingolani, A	2003	1	J Organomet Chem, su	.
Huq, F	1974  4	411	J Cryst Mol Struct	HCA
Maitlis, P	1981  10	1	Chem Soc Rev	HCA
Pettinari, C	2001  4	1290	Inorg Chem Commun	HCA
Pettinari, C	2002  651	5	J Organomet Chem	HCA
Pignolet, L	1983	1	Homogeneous Catalysi	1

Trzeciak, A | 1999 | 190-1 | 883 | Coord Chem Rev | HCA | Westcott, S | 1992 | 114 | 8863 | J Am Chem Soc | HCA

OS.CITING REF COUNT: 5 THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD (5 CITINGS)

L20 ANSWER 3 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 140:156103 HCA Full-text

TITLE: Reactivity of rhodium- $\beta$ -diketonato cyclooctadiene derivatives with mono- and di-phosphines. Synthesis, structural and spectroscopic characterization of Rh(I)

and Rh(III) species containing unsymmetrical

 $\beta$ -diketonate and P-donor ligands

AUTHOR(S): Pettinari, Claudio; Marchetti, Fabio; Pettinari,

Riccardo; Pizzabiocca, Adriano; Drozdov, Andrei;

Troyanov, Sergey I.; Vertlib, Vyatcheslav

CORPORATE SOURCE: Dipartimento di Scienze Chimiche, Universita degli

Studi, Camerino, 62032, Italy

SOURCE: Journal of Organometallic Chemistry (2003),

688(1-2), 216-226

CODEN: JORCAI; ISSN: 0022-328X

PUBLISHER: Elsevier Science B.V.

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 140:156103

AB From the reaction of [Rh(Q)(1,5-cod)] (HQ = 1-phenyl-3-methyl-4-R-pyrazol-5-one: R = 2-thenoyl (HQS) or 2-furanoyl (HQO)) with PPh3, 1,2-bis(diphenylphosphino)ethane (dppe) or 1,3-bis(diphenylphosphino)propane (dppp) in anhydrous solvents under N2, [Rh(Q)(PPh3)2], [Rh(dppe)2](Q) (Q = QS or QO), and [Rh(dppp)(QO)] were obtained. The reactions of [Rh(QS)(1,5-cod)] with CH3I, I2, HCl and C3H5Br in the presence of PPh3 were also studied. All compds. obtained were characterized by elemental analyses, FTIR, ESI-MS spectroscopy, 1H-, 31P- and in selected cases by 13C-NMR spectroscopy. [Rh(QS)(PPh3)2], [Rh(dppe)2](QS), [Rh(QO)(dppp)] and [Rh(QS)C12(PPh3)2], were also characterized in the solid state by single crystal x-ray diffraction. In the air oxidation of [Rh(Q)(PR3)2] and [Rh(QO)(dppp)] occurred, species containing a η2-peroxo group being always identified.

IT 651301-17-6P 651301-21-2P 651301-38-1P 651301-40-5P

(preparation and crystal structure of)

RN 651301-17-6 HCA

CN Rhodium, [2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl- $\kappa$ O)-3H-pyrazol-3-onato- $\kappa$ O3]bis(triphenylphosphine)-, (SP-4-3)- (CA INDEX NAME)

RN 651301-21-2 HCA

CN Rhodium(1+), bis[1,2-ethanediylbis[diphenylphosphine- $\kappa$ P]]-, (SP-4-1)-, salt with 2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl)-

3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 429655-52-7 CMF C15 H11 N2 O2 S

CM 2

CRN 47895-57-8 CMF C52 H48 P4 Rh CCI CCS

RN 651301-38-1 HCA

CN Rhodium, dichloro[2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl- $\kappa$ O)-3H-pyrazol-3-onato- $\kappa$ O3]bis(triphenylphosphine)-, (OC-6-43)- (CA INDEX NAME)

RN 651301-40-5 HCA

CN Rhodium, [4-(2-furanylcarbonyl- $\kappa$ O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ O3][1,3-propanediylbis[diphenylphosphine- $\kappa$ P]]-, (SP-4-3)-, compd. with benzene (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 651301-28-9 CMF C42 H37 N2 O3 P2 Rh CCI CCS

Ph Ph O O O N Ph

CM 2
CRN 71-43-2

C6 H6

CMF

IT 444772-13-8P 444772-14-9P (preparation and reaction with phosphines)

RN 444772-13-8 HCA

CN Rhodium, [(1,2,5,6- $\eta$ )-1,5-cyclooctadiene][2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl- $\kappa$ 0)-3H-pyrazol-3-onato- $\kappa$ 03]- (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

RN 444772-14-9 HCA

CN Rhodium, [(1,2,5,6- $\eta$ )-1,5-cyclooctadiene][4-(2-furanylcarbonyl- $\kappa$ O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ O3]- (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

IT 651301-18-7F 651301-24-5F 651301-28-9F 651301-32-5F 651301-36-9F (preparation of)

RN 651301-18-7 HCA

CN Rhodium, [4-(2-furanylcarbonyl- $\kappa$ 0)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03]bis(triphenylphosphine)-, (SP-4-3)- (CA INDEX NAME)

RN 651301-24-5 HCA

CN Rhodium(1+), bis[1,2-ethanediylbis[diphenylphosphine- $\kappa$ P]]-, (SP-4-1)-, salt with 4-(2-furanylcarbonyl)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 429655-49-2 CMF C15 H11 N2 O3

CM 2

CRN 47895-57-8 CMF C52 H48 P4 Rh CCI CCS

RN 651301-28-9 HCA

CN Rhodium, [4-(2-furanylcarbonyl- $\kappa$ 0)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03][1,3-propanediylbis[diphenylphosphine- $\kappa$ P]]-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 651301-32-5 HCA

CN Rhodium, [2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl-κ0)-3Hpyrazol-3-onato-κ03]diiodobis(triphenylphosphine)-, (OC-6-43)- (CA
INDEX NAME)

$$\begin{array}{c} \text{Ph}_{3P} \\ \text{Ph}_{3P} \\ \text{Ph}_{3P} \\ \text{Ph}_{3P} \\ \text{Ph}_{3P} \\ \text{Ph} \\ \end{array}$$

RN 651301-36-9 HCA

CN Rhodium, bromo[2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl- $\kappa$ O)-3H-pyrazol-3-onato- $\kappa$ O3]( $\eta$ 3-2-propenyl)(triphenylphosphine)- (CA INDEX NAME)

CC 78-7 (Inorganic Chemicals and Reactions)

Section cross-reference(s): 29, 75

IT 651301-17-6P 651301-21-2P 651301-38-1P 651301-40-5P

(preparation and crystal structure of)

IT 444772-13-8P 444772-14-9P

(preparation and reaction with phosphines)

IT 651301-18-7P 651301-24-5P 651301-28-9P

651301-32-5P 651301-36-9P

(preparation of)

RETABLE

Referenced Author	Year   VOL	PG	Referenced Work   Referenced
(RAU)	(RPY) (RVL	)   (RPG)	(RWK)   File
	=+=====+====	=+=====	-+
Anderson, P	1981  20	4101	Inorg Chem
Angermund, K	1997  3	1755	Chem Eur J   HCA
An <b>o</b> n	1995  12	1	Comprehensive Organo
Augustine, R	1970	497	J Chem Soc Chem Comm
Baker, R	1995  34	1336	Angew Chem Int Ed En HCA
Bennett, M	1977  16	1581	Inorg Chem   HCA
Brown, J	1993	125	Chem Soc Rev   HCA
Chin, C	1992	1323	J Chem Soc Dalton Tr HCA
Cingolani, A	2002  41	1151	Inorg Chem   HCA
Cingolani, A	2002  329	100	Inorg Chim Acta   HCA

Crabtree, R	1994	ı	1	The Organometallic C	1
Dai, C	1998			J Chem Soc Chem Conm	
Duan, Z					HCA
	1974			J Chem Soc Dalton Tr	ĺ
Elduque, A	1996			J Chem Soc Dalton Tr	
Esteruelas, M					HCA
Fernandez, E	1997	-		J Chem Soc Chem Comm	HCA
Fornica, R	11995			J Chem Soc Chem Comm	
Fornika, R	2001	79	1642	Can J Chem	HCA
Fryzuk, M			2134		HCA
Giordano, C			188	Inorg Synth	ĺ
Haarman, H	1997				HCA
Hutschka, F					HCA
Jensen, B					HCA
Jesse, A					HCA
John, K				_	HCA
John, K				Organometallics	I
				Rec Trav Chim Pays-b	HCA
Lahuerta, P	-				HCA
Lange, S				J Chem Soc Dalton Tr	•
Leipoldt, J					HCA
Leipoldt, J					HCA
Leitner, W				_	HCA
Marder, T				_	HCA
					HCA
Okafor, E				2	HCA
Pettinari, C					HCA
Pettinari, C				_	HCA
Pettinari, C	-			_	HCA
Ramsden, J	12002	-		J Chem Soc Chem Comm	
Roucoux, A		-			HCA
Senko, M	1 1 2 2 0	1 1 2	•	IsoPro Isotopic Abun	•
Sheldrick, G	  1997	1 1		SHELX-93	1
Sheldrick, G	11997	I I		SHELXS-97	1
Shestakova, E		1   7	-	Rhodium Express	l 1
Shestakova, E					HCA
Shestakova, E				Rhodium Express	I IICA
•	-	-		_	I   HCA
Shestakova, E	1995  2000	-		_	HCA
Simanko, W Suzuki, H	-	-		3	•
	11982				HCA
Togni, A	11994				HCA
Trzeciak, A	11997			J Chem Soc Dalton Tr	
Trzeciak, A	-				HCA
				_	HCA
van Dam, H				_	HCA
van Haaren, R					HCA
van Rooy, A	-		134		HCA
van Rooy, A	-				HCA
Van Vugt, B	-	-	=		HCA
Volger, H	-			_	HCA
Westcott, S					HCA
Westcott, S				_	HCA
					HCA
Zhou, Z				• •	HCA
OS.CITING REF COUNT:	13			3 CAPLUS RECORDS THAT	CITE THIS
		RECO:	RD (13	CITINGS)	

L20 ANSWER 4 OF 16 HCA COPYRIGHT 2011 ACS on STN

ACCESSION NUMBER: 137:263145 HCA Full-text

The reactivity of pay (1.5 cycleses

TITLE: The reactivity of new (1,5-cyclooctadiene)rhodium

acylpyrazolonates towards N- and P-donor ligands:

X-ray structures of [Rh(1, 5-COD)Qs],  $[Rh(1, 5-COD) (phen)]Qs \cdot 0.5H2O (HQs =$ 

1-phenyl-3-methyl-4-(2-thenoyl)-pyrazol-5-one) and

[Rh(1,5-COD)Br]2

AUTHOR(S): Pettinari, Claudio; Marchetti, Fabio; Cingolani,

Augusto; Bianchini, Gianluca; Drozdov, Andrei;

Vertlib, Vyacheslav; Troyanov, Sergei

CORPORATE SOURCE: Dipartimento di Scienze Chimiche, Universita degli

Studi, Camerino MC, 62032, Italy

SOURCE: Journal of Organometallic Chemistry (2002),

651(1-2), 5-14

CODEN: JORCAI; ISSN: 0022-328X

PUBLISHER: Elsevier Science B.V.

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 137:263145

AΒ [Rh(1,5-COD)(Q)] were prepared by the reaction between [Rh(1,5-COD)C1]2 (1,5-COD = 1,5-cyclooctadiene) and HQ (HQ = 1-phenyl-3-methyl-4-R-pyrazol-5-one: R = 2-thenoyl (HQs), 2-furanoyl (HQo) or tert-butylacetyl (HQT)). [Rh(1,5-COD)(Q)] reacted with N2-donor ligands such as 1,10-phenanthroline (phen) or 2,2-bipyridyl (bipy) yielding ionic compds. [Rh(1,5-COD)(N2-donor)]Q. substitutional lability of 1,5-COD in [Rh(1,5-COD)(Q)] vs. mono- and diorganophosphine ligands was also studied. In all cases 1,5-COD was displaced. Reaction with two equivalent of PPh3 gave, upon oxidation of the Rh(I) center, [Rh(PPh3)2(O2)(Q)] species containing a  $\eta2$ -peroxo-group. Reaction of [Rh(1,5-COD)(Q)] with the chelating P2-donor 1,2-bis(diphenylphosphino)ethane (dppe) or 4,4'-bis(diphenylphosphino)ferrocene (dppf) yielded the peroxo Rh(III) compds. [Rh(dppe)202]QT and [Rh(dppf)02(Qs)] or Rh(I) species [Rh(dppf-O2)(QT)] containing the diphosphine in the oxidized form. Finally the reaction between [Rh(1, 5-COD)(Q)] and allyl bromide yielded the known [Rh(1, 5-COD)(Q)]COD)Br]2. All complexes were characterized by anal. and spectral data (IR, 1H and  $31P\{1H\}-NMR$  spectra). The crystal structures of [Rh(1,5-COD)(Qs)], [Rh(1,5-COD)(phen)]Qs and [Rh(1,5-COD)Br]2, all containing a Rh(I) atom in a square coordinate environment, were also reported.

IT 463975-75-9P

(preparation and crystal structure of)

RN 463975-75-9 HCA

CN Rhodium(1+), [(1,2,5,6- $\eta$ )-1,5-cyclooctadiene](1,10-phenanthroline-  $\kappa$ N1, $\kappa$ N10)-, salt with 2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl)-3H-pyrazol-3-one, hydrate (2:2:1) (9CI) (CA INDEX NAME)

CM 1

CRN 463975-74-8

CMF C20 H20 N2 Rh . C15 H11 N2 O2 S

CM 2

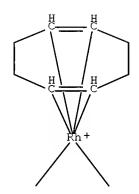
CRN 429655-52-7 CMF C15 H11 N2 O2 S

CM 3

CRN 47248-45-3 CMF C20 H20 N2 Rh

CCI CCS

PAGE 1-A



N N

PAGE 2-A

IT 463975-76-0P 463975-77-1P 463975-78-2P 463975-79-3P 463975-80-6P 463975-81-7P 463975-82-8P 463975-83-9P 463975-84-0P (preparation of)

RN 463975-76-0 HCA

CN Rhodium(1+), [(1,2,5,6- $\eta$ )-1,5-cyclooctadiene](1,10-phenanthroline- $\kappa$ N1, $\kappa$ N10)-, salt with 4-(2-furanylcarbonyl)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 429655-49-2

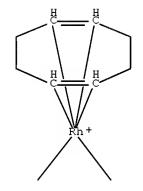
CMF C15 H11 N2 O3

$$\begin{array}{c} \begin{array}{c} \begin{array}{c} \\ \\ \end{array} \end{array} \begin{array}{c} \begin{array}{c} \\ \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \\ \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array} \end{array} \begin{array}{c} \\ \end{array} \begin{array}{c} \\ \end{array}$$

CM 2

CRN 47248-45-3 CMF C20 H20 N2 Rh CCI CCS

PAGE 1-A



PAGE 2-A

RN 463975-77-1 HCA

CN Rhodium(1+), (2,2'-bipyridine- $\kappa$ N1, $\kappa$ N1')[(1,2,5,6- $\eta$ )-1,5-cyclooctadiene]-, salt with 2,4-dihydro-5-methyl-2-phenyl-4-(2-

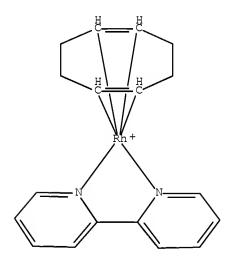
thienylcarbonyl)-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 429655-52-7 CMF C15 H11 N2 O2 S

CM 2

CRN 47101-12-2 CMF C18 H20 N2 Rh CCI CCS



RN 463975-78-2 HCA

CN Rhodium(1+), (2,2'-bipyridine- $\kappa$ N1,  $\kappa$ N1')[(1,2,5,6- $\eta$ )-1,5-cyclooctadiene]-, salt with 4-(2-furanylcarbonyl)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 429655-49-2 CMF C15 H11 N2 O3

CM 2

CRN 47101-12-2 CMF C18 H20 N2 Rh CCI CCS

RN 463975-79-3 HCA

CN Rhodium, [2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl- $\kappa$ O)-3H-pyrazol-3-onato- $\kappa$ O3]peroxybis(triphenylphosphine)- (9CI) (CA INDEX NAME)

RN 463975-80-6 HCA

CN Rhodium, [4-(2-furanylcarbonyl- $\kappa$ 0)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03]peroxybis(triphenylphosphine)- (9CI) (CA INDEX NAME)

RN 463975-81-7 HCA

CN Rhodium, [4-[3,3-dimethyl-1-(oxo-κO)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-κO3]peroxybis(triphenylphosphine)- (9CI) (CA INDEX NAME)

RN 463975-82-8 HCA

CN Rhodium, [1,1'-bis(diphenylphosphino-κP)ferrocene][2,4-dihydro-5-methyl-2-phenyl-4-(2-thienylcarbonyl-κO)-3H-pyrazol-3-onato-κO3]peroxy- (9CI) (CA INDEX NAME)

RN 463975-83-9 HCA

CN Rhodium(1+), bis[1,2-ethanediylbis[diphenylphosphine-κP]]peroxy-, (OC-6-21)-, salt with 4-(3,3-dimethyl-1-oxobutyl)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 216253-41-7 CMF C16 H19 N2 O2

CM 2

CRN 47898-18-0

CMF C52 H48 O2 P4 Rh

CCI CCS

RN 463975-84-0 HCA

CN Rhodium,  $[1,1'-bis(diphenylphosphinyl-\kappa0)ferrocene]$  [4-[3,3-dimethyl-1-(oxo- $\kappa$ 0)butyl]-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03]-, (SP-4-3)- (9CI) (CA INDEX NAME)

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ΙT
     444772-14-9P 463975-73-7P
        (preparation, coordinative substitution reaction with nitrogen and
        phosphorus donor ligands, and oxidation in presence of phosphines)
     444772-14-9 HCA
RN
     Rhodium, [(1,2,5,6-\eta)-1,5-\text{cyclooctadiene}][4-(2-\text{furanylcarbonyl-}
CN
     \kappa0)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa03]-
     (CA INDEX NAME)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
     463975-73-7 HCA
RN
     Rhodium, [(1,2,5,6-\eta)-1,5-\text{cyclooctadiene}][4-[3,3-\text{dimethyl}-1-(\text{oxo-}
CN
     κΟ) butyl]-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-
     κO3]- (CA INDEX NAME)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
     444772-13-8P
TΤ
        (preparation, coordinative substitution reaction with nitrogen and
        phosphorus donor ligands, oxidation in presence of phosphines, and crystal
        structure of)
     444772-13-8 HCA
RN
     Rhodium, [(1,2,5,6-\eta)-1,5-\text{cyclooctadiene}][2,4-\text{dihydro}-5-\text{methyl}-2-
CN
     phenyl-4-(2-thienylcarbonyl-\kappa0)-3H-pyrazol-3-onato-\kappa03]- (CA
     INDEX NAME)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
     29-13 (Organometallic and Organometalloidal Compounds)
     Section cross-reference(s): 75, 78
     12092-45-4P, Bis[bromo(1,5-cyclooctadiene)rhodium] 463975-75-9P
ΤТ
        (preparation and crystal structure of)
     463975-76-0P 463975-77-1P 463975-78-2P
ΙT
     463975-79-3P 463975-80-6P 463975-81-7P
     463975-82-8P 463975-83-9P 463975-84-0P
        (preparation of)
     444772-14-9P 463975-73-7P
ΤТ
        (preparation, coordinative substitution reaction with nitrogen and
        phosphorus donor ligands, and oxidation in presence of phosphines)
ΙT
     444772-13-82
        (preparation, coordinative substitution reaction with nitrogen and
        phosphorus donor ligands, oxidation in presence of phosphines, and crystal
        structure of)
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OS.CITING REF COUNT:
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                              THERE ARE 24 CAPLUS RECORDS THAT CITE THIS
                              RECORD (24 CITINGS)
L20 ANSWER 5 OF 16 HCA COPYRIGHT 2011 ACS on STN
ACCESSION NUMBER: 137:149282 HCA Full-text
                        Complexes of some d and f elements with new
TITLE:
                        4-acylpyrazol-5-ones: synthesis and study
                        Drozdov, A. A.; Vertlib, V. A.; Timokhin, I.;
AUTHOR(S):
                        Troyanov, S. I.; Pettinari, C.; Marchetti, F.
                        Moscow State University, Moscow, 117234, Russia
CORPORATE SOURCE:
SOURCE:
                        Russian Journal of Coordination Chemistry (Translation
                        of Koordinatsionnaya Khimiya) (2002), 28(4),
                        259-263
                        CODEN: RJCCEY; ISSN: 1070-3284
PUBLISHER:
                        MAIK Nauka/Interperiodica Publishing
DOCUMENT TYPE:
                        Journal
LANGUAGE:
                        English
OTHER SOURCE(S):
                        CASREACT 137:149282
     Complexes of 1-phenyl-3-methyl-4-thenoylpyrazol-5-one and 1-phenyl-3-methyl-4-
     furancarbonylpyrazol-5-one, which was synthesized for the 1st time, with Cu,
     Rh, La, and Eu were studied. The substances obtained were studied using
     elemental anal. and IR spectroscopy, while the Rh derivs. were addnl. studied
     using the 1H NMR method. The thermal stability of the Cu derivs. was studied upon heating in a vacuum. The presence of addnl. heteroatoms in
     acylpyrazolone had virtually no effect on the structure and composition of the
     complexes formed. These ligands coordinate atoms of d and f metals through O
     atoms similarly to other \beta-diketones, whereas the remaining heteroatoms only
     participated in the formation of a H bonding system with addnl. ligands or
     coordinated solvent mols.
ΙT
    444772-13-8P 444772-14-9P
       (preparation of)
    444772-13-8 HCA
RN
    Rhodium, [(1,2,5,6-\eta)-1,5-\text{cyclooctadiene}][2,4-\text{dihydro}-5-\text{methyl}-2-
CN
    phenyl-4-(2-thienylcarbonyl-κ0)-3H-pyrazol-3-onato-κ03]- (CA
    INDEX NAME)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
RN
    444772-14-9 HCA
    Rhodium, [(1,2,5,6-\eta)-1,5-\text{cyclooctadiene}][4-(2-\text{furanylcarbonyl-}
CN
    \kappaO)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappaO3]-
    (CA INDEX NAME)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
    78-7 (Inorganic Chemicals and Reactions)
CC
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    444772-10-5P
                  444772-12-7P 444772-13-8P 444772-14-9P
IT
        (preparation of)
RETABLE
  Referenced Author | Year | VOL | PG | Referenced Work | Referenced
     (RAU) | (RPY) | (RVL) | (RPG) | (RWK) | File
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OS.CITING REF COUNT:
                         4
                                (4 CITINGS)
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L20 ANSWER 6 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 135:195655 HCA Full-text

TITLE: Interaction of Rh(I) with a new polydentate O4, N-donor

pyrazolone able to form mononuclear, dinuclear and

heterobimetallic compounds

AUTHOR(S): Pettinari, C.; Marchetti, F.; Drozdov, A.; Vertlib,

V.; Troyanov, S.

CORPORATE SOURCE: Dipartimento di Scienze Chimiche, Universita degli

Studi di Camerino, Camerino, MC, 62032, Italy

SOURCE: Inorganic Chemistry Communications (2001),

4(6), 290-293

CODEN: ICCOFP; ISSN: 1387-7003

PUBLISHER: Elsevier Science S.A.

DOCUMENT TYPE: Journal LANGUAGE: English

OTHER SOURCE(S): CASREACT 135:195655

- AB Reaction of the new polydentate O4,N-donor ligand, namely 2,6-bis[4(1-N-phenyl-3-methyl-pyrazolium-5-one)carbonyl]pyridinium trichloride [H5QN]Cl3, with [Rh(COD)Cl]2 affords the dinuclear compound [Rh2(COD)2([H3QN]Cl3)]1 which reacts with PPh3 forming [Rh(PPh3)2Cl2(HQN)]2, able to interact with SnMe2Br2 yielding the heterobimetallic adduct [Rh(PPh3)2 Cl2(HQN)SnMe2Br2]3; the X-ray structure of 2 shows the Rh(III) center in a slightly distorted octahedral environment with Cl atoms situated in cis-position and two phosphine groups in trans.
- IT 357334-17-9P

(preparation and complexation with tin bromo Me complex)

- RN 357334-17-9 HCA
- CN Rhodium, dichloro[2,4-dihydro-4-[[6-[(5-hydroxy-3-methyl-1-phenyl-1H-pyrazol-4-yl)carbonyl]-2-pyridinyl]carbonyl- $\kappa$ 0]-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03]bis(triphenylphosphine)-, (OC-6-42)- (CA INDEX NAME)

IT 357334-19-1P

(preparation and crystal structure of)

RN 357334-19-1 HCA

CN Rhodium, dichloro[2,4-dihydro-4-[[6-[(5-hydroxy-3-methyl-1-phenyl-1H-pyrazol-4-yl)carbonyl]-2-pyridinyl]carbonyl- $\kappa$ 0]-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03]bis(triphenylphosphine)-, (OC-6-42)-, compd. with 1,1'-oxybis[ethane] (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 357334-17-9

CMF C63 H50 C12 N5 O4 P2 Rh

CCI CCS

CM 2

CRN 60-29-7

CMF C4 H10 O

## IT 357334-16-8P

(preparation and substitution/oxidative addition reaction with chloroform solution

of triphenylphosphine)

RN 357334-16-8 HCA

CN Rhodium, bis[ $(1,2,5,6-\eta)-1$ ,5-cyclooctadiene][ $\mu$ -[[4,4'-[2,6-pyridinediyldi(carbonyl- $\kappa$ 0)]bis[2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03]](2-)]]di-, trihydrochloride (9CI) (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

IT 357334-18-0P

(preparation of)

RN 357334-18-0 HCA

CN Rhodium, dichloro(dibromodimethyltin)[ $\mu$ -[[4,4'-[2,6-pyridinediyldi(carbonyl- $\kappa$ O)]bis[2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ O3]](2-)]]bis(triphenylphosphine)-, stereoisomer (9CI) (CA INDEX NAME)

CC 29-13 (Organometallic and Organometalloidal Compounds) Section cross-reference(s): 75

IT 357334-17-9P

(preparation and complexation with tin bromo Me complex)

IT 357334-19-1P

(preparation and crystal structure of)

IT 357334-16-8P

 $\hbox{ (preparation and substitution/oxidative addition reaction with chloroform solution }$ 

of triphenylphosphine)

IT 357334-18-0P

(preparation of)

RETABLE

KETVDIE				
Referenced Author	Year   VOL	PG	Referenced Work	Referenced
(RAU)	(RPY) (RVL)	(RPG)	(RWK)	File
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OS.CITING REF COUNT:	7	THE	RE ARE	7 CAPLUS RECORDS THAT	CITE THIS RECORD			
(7 CITINGS)								

L20 ANSWER 7 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 132:259719 HCA Full-text

TITLE: Synthesis and cytotoxic activity of some novel

trans-palladium complexes with pyrazole derivatives

AUTHOR(S): Al-Allaf, Talal A. K.; Rashan, Luay J.

CORPORATE SOURCE: Department of Chemistry, College of Science, Applied

Science University, Amman, 11931, Jordan Asian Journal of Chemistry (1999), 11(4),

1543-1545

CODEN: AJCHEW; ISSN: 0970-7077

PUBLISHER: Asian Journal of Chemistry

DOCUMENT TYPE: Journal LANGUAGE: English

AB Novel Pd(II) complexes trans-[PdL2C12], where L is a pyrazole derivative, were prepared and characterized physicochem. and spectroscopically. Pyrazole derivs. coordinate with Pd in a monodentate fashion via the most reactive N site. The cytotoxic activity of these complexes was evaluated in vitro against four cell-lines using the MTT-assay, one fluid suspension (P388, leukemia) and three solid human cell lines (Hep-2, larynx; RD, embryonal rhabdomyosarcoma and HeLa, cervical cells). One of these complexes, for example, demonstrated a potent cytotoxic activity against P388 and significant cytotoxicity against the other three-cell lines in comparison with the reference stds.: cisplatin, carboplatin, oxaliplatin and 5-FU.

IT 262597-31-9P

SOURCE:

(preparation and antitumor activity)

RN 262597-31-9 HCA

CN Palladium, dichlorobis(3,6-dimethylpyrano[2,3-c]pyrazol-4(1H)-one- $\kappa$ N2)-, (SP-4-1)- (CA INDEX NAME)

$$\begin{array}{c|c} & & & & \\ & &$$

Section cross-reference(s): 1

IT 121833-06-5P 126873-41-4P 262597-25-1P 262597-27-3P 262597-29-5P 262597-30-8P 262597-33-9P 262597-32-0P 262597-33-1P

(preparation and antitumor activity)

RETABLE

Referenced Author	Year	VOL	PG	Referenced Work	Referenced		
(RAU)	(RPY)	(RVL)	)   (RPG)	(RWK)	File		
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OS.CITING REF COUNT:	3	THE	RE ARE	3 CAPLUS RECORDS THAT	CITE THIS RECORD		
(3 CITINGS)							

L20 ANSWER 8 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 129:316354 HCA Full-text

ORIGINAL REFERENCE NO.: 129:64559a,64562a

TITLE: (1-Phenyl-3-methyl-4-acetylpyrazolon-5-ato)rhodium(I)

complexes, synthesis, structural and spectroscopical

characterization: Reactivity of diolefin- and dicarbonyl-rhodium complexes toward N-, P- and

0-donors

AUTHOR(S): Pettinari, C.; Accorroni, F.; Cingolani, A.;

Marchetti, F.; Cassetta, A.; Barba, L.

CORPORATE SOURCE: Dipartimento di Scienze Chimiche, Universita di

Camerino, Camerino, I-62032, Italy

SOURCE: Journal of Organometallic Chemistry (1998),

566(1-2), 187-201

CODEN: JORCAI; ISSN: 0022-328X

PUBLISHER: Elsevier Science S.A.

DOCUMENT TYPE: Journal LANGUAGE: English

Novel complexes of rhodium(I) [Rh(diolefin)(Q'')] (where HQ'' = 1-phenyl-3methyl-4-acetylpyrazol-5-one and diolefin = cycloocta-1,5-diene (COD), bicyclo[2.2.1]hepta-2,5-diene (NBD) or 1,5-hexadiene (HEX)) were synthesized and characterized by anal. and spectral data. [Rh(COD)(Q'')] interacts with 4,5-dimethyl-1,10-phenanthroline (Me2Phen) and 2,2'-bipyridyl (Bipy) yielding the cationic derivs. [Rh(COD)(Me2Phen)](Q'')(H2O), [Rh(COD)(Bipy)](Q'')(H2O) upon displacement of the (Q'') - donor from the coordination sphere of the metal center. Whereas [Rh(COD)(Q'')] interacts with 2-benzoylpyridine (Bzpy) yielding the 1:1 adduct [Rh(COD)(Bzpy)(Q'')] in which Bzpy acts as Nmonodentate donor. On the other hand the monodentate P-donors triphenylphosphine, tri-Ph phosphite, tricyclohexylphosphine and the bidentate bis(diphenylphosphino)ethane (DPPE) displace the COD ligand from [Rh(COD)(Q'')] giving the neutral derivs. [Rh(PR3)2(Q'')] (PR3 = PPh3, or P(OPh)3) and [Rh(DPPE)(Q'')](H2O). HQ'' reacts with the dinuclear [Rh(CO)2Cl]2. The tetradentate cyclooctatetraene (COT) reacts with [Rh(CO)2(Q'')] yielding the derivative [Rh(CO)2(HQ'')Cl] in which HQ'' acts as neutral monodentate O-donor ligand. Whereas in presence of NEt3, HQ'' reacts

with [Rh(CO)2C1]2 yielding [Rh(CO)2(Q'')]. In this complex, one mol. of CO can be replaced by one mole of Phen and Bipy or by two moles of PPh3 and AsPh3 yielding the derivs.  $[Rh(CO)(L)n(Q'')] \cdot x(H2O)$  (L = Me2Phen or Bipy, n = 1; L = PPh3 or AsPh3, n = 2) whereas one mole of DPPE displaces both the mols. of CO, yielding [Rh(DPPE)(Q'')] yielding the derivative [Rh(COT)(Q'')]. The x-ray crystal structure determination of [Rh(COD)(Q'')] establishes that the rhodium atom is in a square planar configuration with two adjacent sites occupied by the (Q'')- ligand in the O2-bidentate form (Rh-O) distances = 2.054(2) and 2.061(2) A). The COD ring has a twisted boat conformation with Rh-C distances in the range 2.101(3)-2.110(3) A. Comparison was made with structural data reported for several related tetracoordinated (COD)Rh(I) adducts.

IT 214747-44-1P

(crystal structure; preparation, structural, and spectroscopical characterization of acetylpyrazolonato rhodium complex and reactivity of diolefin- and dicarbonyl-rhodium complexes toward nitrogen-, phosphorus- and oxygen-donors)

RN 214747-44-1 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3][(1,2,5,6-<math>\eta$ )-1,5-cyclooctadiene]- (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

IT 214747-45-2P 214747-46-3P 214747-48-5P 214747-49-6P 214747-50-9P 214747-51-0P 214747-52-1P 214747-53-2P 214747-54-3P 214747-56-5P 214747-57-6P 214747-58-7P 214747-59-8P 214747-60-1P

(preparation of)

RN 214747-45-2 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3][(1,2,5,6-\eta)-1,5-hexadiene]- (CA INDEX NAME)$ 

$$\begin{array}{c} & & \\$$

RN 214747-46-3 HCA

CN Rhodium, [4-(acetyl- $\kappa$ 0)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03][(2,3,5,6- $\eta$ )-bicyclo[2.2.1]hepta-2,5-diene]- (CA INDEX NAME)

RN 214747-48-5 HCA

CN Rhodium(1+), [(1,2,5,6- $\eta$ )-1,5-cyclooctadiene](4,7-dimethyl-1,10-phenanthroline- $\kappa$ N1, $\kappa$ N10)-, salt with 4-acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

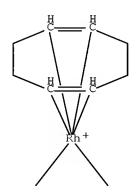
CM 1

CRN 214747-47-4 CMF C12 H11 N2 O2

CM 2

CRN 56678-53-6 CMF C22 H24 N2 Rh CCI CCS

PAGE 1-A



PAGE 2-A

RN 214747-49-6 HCA

CN Rhodium(1+), (2,2'-bipyridine- $\kappa$ N1, $\kappa$ N1')[(1,2,5,6- $\eta$ )-1,5-cyclooctadiene]-, salt with 4-acetyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 214747-47-4 CMF C12 H11 N2 O2

CM 2

CRN 47101-12-2 CMF C18 H20 N2 Rh

CCI CCS

RN 214747-50-9 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3][(1,2,5,6-\eta)-1,5-cyclooctadiene][phenyl(2-pyridinyl-$ 

 $\kappa$ N)methanone]- (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

RN 214747-51-0 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3]bis(triphenylphosphine)-, (SP-4-3)- (CA INDEX NAME)$ 

RN 214747-52-1 HCA

CN Rhodium, [ $4-(acetyl-\kappa O)-2$ ,  $4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3]bis(triphenyl phosphite-<math>\kappa P$ )-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 214747-53-2 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa O3$ ] [1,2-ethanediylbis[diphenylphosphine- $\kappa P$ ]]-, (SP-4-3)-(9CI) (CA INDEX NAME)

RN 214747-54-3 HCA

CN Rhodium, dicarbonylchloro[1-[5-(hydroxy- $\kappa$ 0)-3-methyl-1-phenyl-1H-pyrazol-4-yl]ethanone]-, (SP-4-3)- (CA INDEX NAME)

RN 214747-56-5 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3]carbonyl(4,7-dimethyl-1,10-phenanthroline- <math>\kappa N1,\kappa N10)-$ , (SP-5-43)- (CA INDEX NAME)

RN 214747-57-6 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-<math>\kappa O3]$  (2,2'-bipyridine- $\kappa N1,\kappa N1'$ ) carbonyl-, (SP-5-43)- (CA INDEX NAME)

RN 214747-58-7 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3]$  carbonylbis(triphenylphosphine)-, (SP-5-43)- (CA INDEX NAME)

RN 214747-59-8 HCA

CN Rhodium, [4-(acetyl-κ0)-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-κ03]carbonylbis(triphenylarsine)-, (SP-5-43)- (CA INDEX NAME)

RN 214747-60-1 HCA

CN Rhodium, [4-(acetyl- $\kappa$ 0)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato- $\kappa$ 03][(1,2,5,6- $\eta$ )-1,3,5,7-cyclooctatetraene]- (9CI) (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

IT 214747-55-4P

(preparation, structural, and spectroscopical characterization of acetylpyrazolonato rhodium complex and reactivity of diolefin- and dicarbonyl-rhodium complexes toward nitrogen-, phosphorus- and oxygen-donors)

RN 214747-55-4 HCA

CN Rhodium,  $[4-(acetyl-\kappa O)-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-\kappa O3]dicarbonyl-, (SP-4-3)- (CA INDEX NAME)$ 

CC 29-13 (Organometallic and Organometalloidal Compounds)
Section cross-reference(s): 75

IT 214747-44-1P

(crystal structure; preparation, structural, and spectroscopical characterization of acetylpyrazolonato rhodium complex and reactivity of diolefin- and dicarbonyl-rhodium complexes toward nitrogen-, phosphorus- and oxygen-donors)

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IT 214747-45-2P 214747-46-3P 214747-48-5P 214747-49-6P 214747-50-9P 214747-51-0P 214747-52-1P 214747-53-2P 214747-54-3P 214747-56-5P 214747-57-6P 214747-58-7P 214747-59-8P 214747-60-1P (preparation of)
IT 214747-55-4P
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(preparation, structural, and spectroscopical characterization of acetylpyrazolonato rhodium complex and reactivity of diolefin- and dicarbonyl-rhodium complexes toward nitrogen-, phosphorus- and oxygen-donors)

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                                THERE ARE 25 CAPLUS RECORDS THAT CITE THIS
OS.CITING REF COUNT:
                       25
                            RECORD (25 CITINGS)
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L20 ANSWER 9 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 129:194249 HCA Full-text
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ORIGINAL REFERENCE NO.: 129:39361a,39364a TITLE: Thermoanalysis of

bis[1-phenyl-3-methyl-4-benzoylpyrazolone-

5]palladium(II)

AUTHOR(S): Tian, Xin; Jiang, Xuchuan; Yang, Yanzhao; Shao, Hua;

Yang, Yonghui; Sun, Sixiu

CORPORATE SOURCE: School Chem., Shandong Univ., Jinan, Peop. Rep. China

SOURCE: Shandong Daxue Xuebao, Ziran Kexueban (1998)

), 33(2), 201-205

CODEN: SDXKEU; ISSN: 0559-7234

PUBLISHER: Shandong Daxue

DOCUMENT TYPE: Journal LANGUAGE: Chinese

AB Synthesis, IR spectra and thermoanal. of bis-(1-phenyl-3-methyl-4-benzoylpyrazolone-5)palladium(II) were reported. Kinetic parameters were obtained from anal. of the TG, DTG curves by integral and differential methods. The possible reaction mechanism was suggested by comparison with the kinetic parameters.

IT 72585-53-69

(thermal anal. of bis[1-phenyl-3-methyl-4-benzoylpyrazolone-5]palladium(II))

RN 72585-53-6 HCA

CN Palladium, bis  $[4-(benzoy1-\kappa O)-2, 4-dihydro-5-methy1-2-pheny1-3H-$ 

pyrazol-3-onato- $\kappa$ O3]- (CA INDEX NAME)

CC 67-3 (Catalysis, Reaction Kinetics, and Inorganic Reaction Mechanisms)

Section cross-reference(s): 69, 78

IT 72585-53-6P

PUBLISHER:

(thermal anal. of bis[1-phenyl-3-methyl-4-benzoylpyrazolone-5]palladium(II))

L20 ANSWER 10 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 123:159338 HCA Full-text

ORIGINAL REFERENCE NO.: 123:28063a,28066a

TITLE: Studies on platinum(II) and palladium(II) complexes of

some substituted pyrazole-5-ones, pyrazoles, (hydroxyaryl)pyrazoles and pyranopyrazole

AUTHOR(S): Al-Allaf, Talal A. K.; Al-Bayati, Redha I. H.

CORPORATE SOURCE: College of Science, University of Mosul, Mosul, Iraq

SOURCE: Asian Journal of Chemistry (1995), 7(3),

465 - 70

CODEN: AJCHEW; ISSN: 0970-7077 Asian Journal of Chemistry

DOCUMENT TYPE: Journal LANGUAGE: English

AB The coordination behavior of several pyrazole-5-ones and pyrazoles derivs. with Pt(II) and Pd(II) metals are reported by the isolation and characterization of the resulting complexes. These complexes possess a square planar structure (cis-form) as revealed from IR and NMR spectral data. The ligands are coordinated mainly through the N-N linkage of the pyrazole ring.

IT 166898-94-8P 166899-03-2P

(preparation of)

RN 166898-94-8 HCA

CN Platinum, dichloro(methyl 2,3-dihydro-3-oxo-1H-pyrazole-4-carboxylate-N1,N2)-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 166899-03-2 HCA

CN Palladium, dichloro(3,6-dimethylpyrano[2,3-c]pyrazol-4(1H)-one-N1,N2)-, (SP-4-3)- (9CI) (CA INDEX NAME)

CC 78-7 (Inorganic Chemicals and Reactions)

IT 166898-93-7P 166898-94-89 166898-95-9P 166898-96-0P

166898-97-1P 166898-98-2P 166898-99-3P 166899-00-9P 166899-01-0P

166899-02-1P 166899-03-2P 166899-04-3P

(preparation of)

OS.CITING REF COUNT: 5 THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD

(5 CITINGS)

L20 ANSWER 11 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 115:222118 HCA Full-text

ORIGINAL REFERENCE NO.: 115:37629a,37632a

TITLE: Polarographic adsorptive wave of

osmium(VIII)-4,4'-decanedioylbis(1-phenyl-3-methyl-5-

pyrazolone) (DBPMP) complex by using 1.5 order

differential technique Lu, Wen; Wang, Zhaizhong

CORPORATE SOURCE: Dep. Chem., Yunnan Univ., Kunming, 650091, Peop. Rep.

China

SOURCE: Fenxi Shiyanshi (1990), 9(2), 1-5

CODEN: FENSE4; ISSN: 1000-0720

DOCUMENT TYPE: Journal LANGUAGE: Chinese

AB In the 0.2 mol/L NaOH solution containing 3.3+10-5 mol/L DBPMP, a sensitive adsorptive wave of Os(VIII)-DBPMP complex was obtained by using single-sweep polarog. The peak potential is -0.79 V(vs. SCE). The peak height of 1.5 order derivative wave is linearly proportional to the Os(VIII) concentration in the range from 1+10-8 to 1+10-5 mol/L. In the solution, the composition of the complex was determined as [OsO4(DBPMP)]2-. The electrode process and mechanism were studied. It shows that the complex is adsorbed onto the DME; the organic moiety of the complex reduces irreversibly, and the reduced product is also adsorbed to the DME.

IT 136901-58-1

AUTHOR(S):

(polarog. of)

RN 136901-58-1 HCA

CN Osmate(2-), [1,10-bis(4,5-dihydro-3-methyl-5-oxo-1-phenyl-1H-pyrazol-4-yl)-1,10-decanedionato(2-)-0,0',0'',0''']tetraoxo-(9CI) (CA INDEX NAME)

CC 79-6 (Inorganic Analytical Chemistry)

Section cross-reference(s): 72

IT 136901-58-1

(polarog. of)

L20 ANSWER 12 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 114:198653 HCA  $\underline{\text{Full-text}}$ 

ORIGINAL REFERENCE NO.: 114:33271a,33274a

TITLE: Study on the polarographic adsorptive complex wave of

rhodium(III)-DBPMP complexone

AUTHOR(S): Lu, Wen; Wang, Zaizhong

CORPORATE SOURCE: Dep. Chem., Yunnan Univ., Kunming, Peop. Rep. China

SOURCE: Guijinshu (1989), 10(3), 31-9 CODEN: GUIJE7; ISSN: 1004-0676

DOCUMENT TYPE: Journal LANGUAGE: Chinese

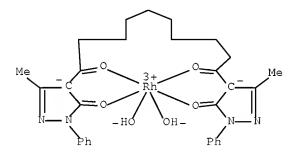
4.4'-Decanedioylbis[1-phenyl-3-methyl-5-pyrazolinone] (DBPMP) in LiCl-NaOH solution produced 2 cathodic reduction waves at -0.43 and -1.28 V, resp. After addition of Rh(III) the wave height of the 2 waves increased with increase of Rh(III) concentration. There was a linearity between the wave height of the first wave with the concentration of Rh(III) in the range 6.0 + 10-9-1.0 + 10-7 mol/L. The UV and IR spectra showed the formation of a Rh-DBPMP complex. The composition of the complex was Rh(III):DBPMP = 1:1 and the stability constant was 1.26 + 106. The effect of surfactants (cetylpyridinium bromide, animal glue, Na dodecanesulfonate) and temperature on the waveheight was studied.

IT 133317-70-1

(IR and UV spectra and stability constant and polarog. of)

RN 133317-70-1 HCA

CN Rhodate(1-), [1,10-bis(4,5-dihydro-3-methyl-5-oxo-1-phenyl-1H-pyrazol-4-yl)-1,10-decanedionato(2-)-0,0',0'',0''']dihydroxy- (9CI) (CA INDEX NAME)



CC 79-6 (Inorganic Analytical Chemistry)

Section cross-reference(s): 68, 72

IT 133317-70-1

(IR and UV spectra and stability constant and polarog. of)

L20 ANSWER 13 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 103:123680 HCA  $\underline{Full-text}$ 

ORIGINAL REFERENCE NO.: 103:19789a,19792a

TITLE: Rhodium and iridium complexes containing the anion of

1-phenyl-3-methyl-4-benzoyl-5-pyrazolone, a

sophisticated analog of  $\beta$ -diketones

AUTHOR(S): Bonati, Flavio; Oro, Luis A.; Pinillos, M. Teresa

CORPORATE SOURCE: Dip. Sci. Chim., Univ. Camerino, Camerino, 62032,

Italy

SOURCE: Polyhedron (1985), 4(2), 357-64 CODEN: PLYHDE; ISSN: 0277-5387

DOCUMENT TYPE: Journal LANGUAGE: English

Several (diolefin)M(A) complexes (M = Rh, Ir; diolefin = e.g. 1,5-cyclooctadiene, 2,5-norbornadiene; AH = 1-phenyl-3-methyl-4-benzoyl-5-pyrazolone, a very stable asym. analog of acetylacetone) were prepared In these complexes the diolefin could be replaced by one mole of (Ph2PCH2CH2)2, two of CO or of PPh3, or three of CNCMe3, whereas 1,10-phenanthroline displaced the chelating ligand to yield [(cyclooctadiene)Rh(phen)]+ (A)-. Some compds. X-Y (X-Y = iodine or MeI) added oxidatively yielding the corresponding trivalent species. Using 31P NMR spectra the presence of the expected steric isomers was detected in (Ph3P)(CO)Rh(A) and in (Ph3P)(CO)Rh(A)(X)(Y).

IT 98063-63-9P 98063-64-0P 98063-65-1P 98063-67-3P 98063-68-4P 98063-69-5P 98063-70-8P 98063-71-9P 98063-72-0P 98063-86-6P 98091-97-5P 98091-98-6P 98091-99-7P 98092-01-4P 98092-03-6P 98092-04-7P 98092-05-8P

98104-43-9P 98111-79-6P

(preparation of)

RN 98063-63-9 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-O,O')carbonyl(triphenylphosphine)-, (SP-4-2)- (9CI) (CA INDEX NAME)

RN 98063-64-0 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')bis(n2-ethene)- (9CI) (CA INDEX NAME)

RN 98063-65-1 HCA

CN Rhodium,  $(4-benzoyl-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')bis[(1,2-<math>\eta$ )-cyclooctene]- (9CI) (CA INDEX NAME)

RN 98063-67-3 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')bis(η3-2-propenyl)- (9CI) (CA INDEX NAME)

$$\begin{array}{c} \text{Ph} \\ \text{N} \\ \text{Me} \end{array}$$

RN 98063-68-4 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')bis(triphenylphosphine)-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 98063-69-5 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[1,4-butanediylbis[diphenylphosphine]-P,P']-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 98063-70-8 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')tris(2-isocyano-2-methylpropane)- (9CI) (CA INDEX NAME)

$$t-Bu-N \stackrel{+}{=} C^ t-Bu-N \stackrel{+}{=} C^ t-Bu-N \stackrel{+}{=} C^ ph$$
 $t-Bu-N \stackrel{+}{=} C^ ph$ 
 $ph$ 
 $ph$ 

RN 98063-71-9 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')dicarbonyl-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 98063-72-0 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')carbonyliodomethyl(triphenylphosphine)- (9CI) (CA INDEX NAME)

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')carbonyldiiodo(triphenylphosphine)- (9CI) (CA INDEX NAME)

RN 98091-97-5 HCA

CN Rhodium,  $(4-benzoyl-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[(1,2,5,6-\eta)-1,5-cyclooctadiene]-(9CI)$  (CA INDEX NAME)

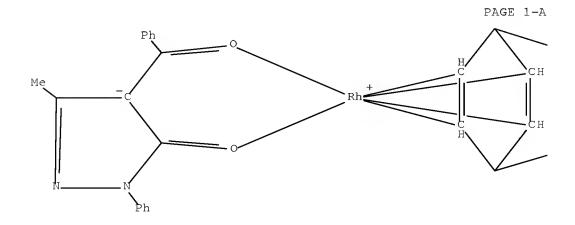
\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

RN 98091-98-6 HCA

CN Rhodium,  $(4-benzoyl-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[(2,3,5,6-\eta)-bicyclo[2.2.1]hepta-2,5-diene]-(9CI)$  (CA INDEX NAME)

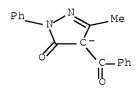
RN 98091-99-7 HCA

CN Rhodium, (4-benzoyl-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0') [(2,3,9,10- $\eta$ )-5,6,7,8-tetrafluoro-1,4-dihydro-1,4-ethenonaphthalene]- (9CI) (CA INDEX NAME)



PAGE 1-B

98092-00-3 HCA RN Rhodium, bis(4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-CN 0,0') [ $\mu$ -[(1,2,5,6- $\eta$ :3,4,7,8- $\eta$ )-1,3,5,7-cyclooctatetraene]]di-(9CI) (CA INDEX NAME) \*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\* 98092-01-4 HCA RN CN Iridium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[(1,2,5,6- $\eta$ )-1,5-cyclooctadiene]- (9CI) (CA INDEX NAME) \*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\* 98092-03-6 HCA RN CN Rhodium(1+),  $(2,2'-bipyridine-N,N')[(1,2,5,6-\eta)-1,5-cyclooctadiene]-$ , salt with 4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-one (1:1) (9CI) (CA INDEX NAME) CM1 CRN 98092-02-5 CMF C17 H13 N2 O2

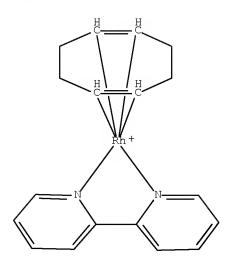


CM 2

CRN 47101-12-2

CMF C18 H20 N2 Rh

CCI CCS



RN 98092-04-7 HCA

CN Rhodium,  $(4-benzoyl-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[(1,2,5,6-\eta)-1,5-cyclooctadiene]diiodo-(9CI)$  (CA INDEX NAME)

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

RN 98092-05-8 HCA

CN Iridium,  $(4-benzoyl-2, 4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[(1,2,5,6-\eta)-1,5-cyclooctadiene]diiodo-(9CI)$  (CA INDEX NAME)

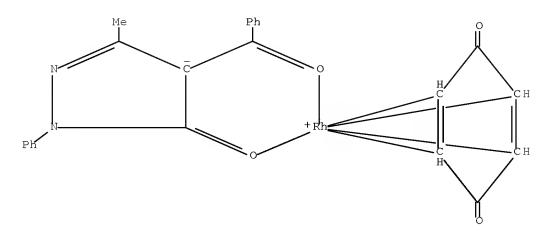
\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

RN 98104-43-9 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')carbonyl(triphenylphosphine)-, (SP-4-3)- (9CI) (CA INDEX NAME)

RN 98111-79-6 HCA

CN Rhodium, (4-benzoyl-2,4-dihydro-5-methyl-2-phenyl-3H-pyrazol-3-onato-0,0')[(2,3,5,6- $\eta$ )-2,5-cyclohexadiene-1,4-dione]- (9CI) (CA INDEX NAME)



CC 29-13 (Organometallic and Organometalloidal Compounds)

IT 98063-63-9P 98063-64-0P 98063-65-1P 98063-66-2P 98063-67-3P 98063-68-4P 98063-69-5P 98063-70-8P 98063-71-9P 98063-72-0P 98063-86-6P 98091-97-5P 98091-98-6P 98091-99-7P 98092-00-3P

98092-01-4P 98092-03-6P 98092-04-7P 98092-05-8P 98104-43-9P 98111-79-6P

(preparation of)

OS.CITING REF COUNT: 14 THERE ARE 14 CAPLUS RECORDS THAT CITE THIS RECORD (14 CITINGS)

L20 ANSWER 14 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 97:109915 HCA Full-text

ORIGINAL REFERENCE NO.: 97:18285a,18288a TITLE: Observations of

1-phenyl-3-methyl-4-trifluoroacetyl-5-pyrazolone. A

promising extracting agent

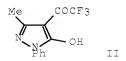
AUTHOR(S): Okafor, Emmanuel Chukwuemeka

CORPORATE SOURCE: Dep. Chem., Univ. Nigeria, Nsukka, Nigeria

SOURCE: Talanta (1982), 29(4), 275-8 CODEN: TLNTA2; ISSN: 0039-9140

DOCUMENT TYPE: Journal LANGUAGE: English

GΙ



4-Trifluoroacetyl-3-methyl-1-phenyl-5-pyrazolone (I), a promising metal extractant, was obtained in 91% yield by treating 1-phenyl-3-methyl-5-pyrazolone with (F3CO)2O in pyridine. Recrystn. studies revealed that only one tautomer, the enol II, can be isolated, sometimes with 1 mol. of water of crystallization, contrary to reports (Jensen, B. S., 1959) that a yellow enol

and a white keto tautomer can be obtained from n-hexane and aqueous EtOH, resp. The m.ps. and colors of some of the metal chelates of I are tabulated and, in the case of Hg(II) and Cu(II) chelates, differ from those reported by others. Solubility data for some of the metal chelates are also given.

ΙT 77259-30-4

(solubility and phys. properties of)

RN 77259-30-4 HCA

Rhodium, tris[2,4-dihydro-5-methyl-2-phenyl-4-(trifluoroacetyl)-3H-pyrazol-CN 3-onato-0,0']- (9CI) (CA INDEX NAME)

28-8 (Heterocyclic Compounds (More Than One Hetero Atom)) CC

Section cross-reference(s): 54

77259-28-0 77259-29-1 **77259-30-4** 77259-31-5 77273-41-7 ΙT 81714-08-1 81714-09-2 81714-06-9 81714-07-0 81714-14-9

81714-15-0 81999-83-9 81999-84-0 81999-88-4

(solubility and phys. properties of)

THERE ARE 5 CAPLUS RECORDS THAT CITE THIS RECORD OS.CITING REF COUNT: 5 (6 CITINGS)

L20 ANSWER 15 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 94:184733 HCA Full-text

ORIGINAL REFERENCE NO.: 94:30087a,30090a

TITLE: The metal complexes of heterocyclic  $\beta$ -diketones

and their derivatives. Part VIII. Synthesis,

structure, proton NMR and infrared spectral studies of

the complexes of aluminum(III), iron(III), cobalt(III), rhodium(III), indium(III), and

zirconium(IV) with

1-phenyl-3-methyl-4-trifluoroacetyl-5-pyrazolone

(HPMTFP)

Okafor, Emmanuel Chukwuemeka AUTHOR(S):

Fac. Phys. Sci., Univ. Nigeria, Nsukka, Nigeria CORPORATE SOURCE:

Zeitschrift fuer Naturforschung, Teil B: Anorganische SOURCE:

Chemie, Organische Chemie (1981), 36B(2),

213 - 17

CODEN: ZNBAD2; ISSN: 0340-5087

DOCUMENT TYPE: Journal LANGUAGE: English

ML3 (M = Al, Fe, Co, Rh, In; HL = 1-phenyl-3-methyl-4-trifluoroacetyl-5pyrazolone) and ZrL4 were prepared and characterized by elemental analyses,

conductivity and magnetic moment measurements, 1H NMR and IR spectroscopy. HLreacts as a bidentate enol forming neutral metal chelates. The 1H NMR spectra of chelates sufficiently soluble in deuterated NMR solvents were recorded and studied. The IR spectra were measured between 4000-200 cm-1 and assignments are proposed for the observed frequencies. The M-O stretching frequency follows the order: Al > Rh > Fe = Co = Zr > In.

IT 77259-30-4P

(preparation of)

RN 77259-30-4 HCA

Rhodium, tris[2,4-dihydro-5-methyl-2-phenyl-4-(trifluoroacetyl)-3H-pyrazol-CN 3-onato-0,0']- (9CI) (CA INDEX NAME)

CC 78-7 (Inorganic Chemicals and Reactions)

77259-28-0P 77259-29-1P **77259-30-4P** 77259-31-5P ΙT

77259-32-6P 77273-41-7P

(preparation of)

THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD OS.CITING REF COUNT:

(2 CITINGS)

L20 ANSWER 16 OF 16 HCA COPYRIGHT 2011 ACS on STN ACCESSION NUMBER: 92:68757 HCA Full-text

ORIGINAL REFERENCE NO.: 92:11199a,11202a

TITLE: Extraction of palladium(II) with

1-phenyl-3-methyl-4-benzoylpyrazole-5-one

Mirza, M. Y.; Bailey, R. T. AUTHOR(S):

Dep. Chem., Univ. Nigeria, Nsukka, Nigeria CORPORATE SOURCE: Journal of Inorganic and Nuclear Chemistry ( SOURCE:

**1979**), 41(5), 772-3

CODEN: JINCAO; ISSN: 0022-1902

DOCUMENT TYPE: Journal LANGUAGE: English

AB Pd pyrazolonate was prepared by extraction of PdC12 from aqueous solution with 1-phenyl-3-methyl-4-benzoylpyrazol-5-one. IR and NMR spectra of the free ligand and Pd pyrazolonate suggested that the ligand coordinates through the 2 O atoms.

72585-53-6P ΙT

(preparation of)

RN 72585-53-6 HCA

CN Palladium, bis[4-(benzoyl-ĸ0)-2,4-dihydro-5-methyl-2-phenyl-3Hpyrazol-3-onato-κ03]- (CA INDEX NAME)

CC 78-7 (Inorganic Chemicals and Reactions)

IT 72585-53-62

(preparation of)

OS.CITING REF COUNT: 8 THERE ARE 8 CAPLUS RECORDS THAT CITE THIS RECORD

(8 CITINGS)